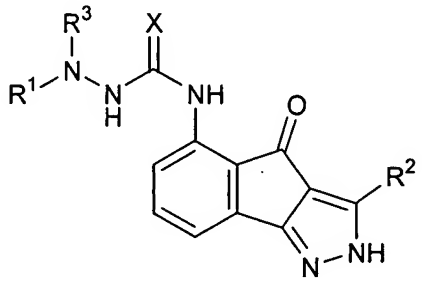
Amendments to the Claims:

This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

Claim 1 (previously presented): A compound according to formula (I):



(I)

X is selected from O or S;

R^1 is selected from the groups: C_3-C_{10} membered carbocycle substituted with 0-5 R^4, and 3-10 membered heterocycle substituted with 0-5 R^5, provided that if R^1 is phenyl then R^1 is substituted with 1-5 R^4;

R^2 is selected from the groups: H, C_{1-10} alkyl substituted with 0-3 R^6, C_{2-10} alkenyl substituted with 0-3 R^6, C_{2-10} alkynyl substituted with 0-3 R^6, $-(CF_2)_mCF_3$, C_{3-10} membered carbocycle substituted with 0-5 R^4, and 3-10 membered heterocycle containing from 1-4 heteroatoms selected from O, N, and S and substituted with 0-5 R^5;

R^3 is selected from the groups: H, C_{1-4} alkyl, C_{3-6} cycloalkyl, or C_{4-10} cycloalkylalkyl;

R^4 is independently selected from the groups: halo, -CN, NO_2, C_{1-4} alkyl, C_{1-4} haloalkyl, NR^7R^{7a}, =O, OR^7, COR^7, CO_2R^7, $CONR^7R^{7a}$, $NHC(O)NR^7R^{7a}$, $NHC(S)NR^7R^{7a}$,

$NR^7C(O)OR^{7b}$, $NR^7C(O)R^{7b}$, $SO_2NR^7R^{7a}$, SO_2R^{7b}, and 5-10 membered heterocycle containing from 1-4 heteroatoms selected from O, N, and S;

alternatively, when two R^4's are present on adjacent carbon atoms they combine to form -OCH_2O- or -OCH_2CH_2O-;

R^5 is independently selected from the groups: halo, -CN, NO_2, C_{1-4} alkyl, C_{1-4} haloalkyl, NR^7R^{7a}, $NR^7C(O)OR^{7b}$, $NR^7C(O)R^{7b}$, OR^7, COR^7, CO_2R^7, $CONR^7R^{7a}$, $CON(R^9)[(CH_2)_mR^{10}]$, $CO(CH_2)_mR^{10}$, $NHC(O)NR^7R^{7a}$, $NHC(S)NR^7R^{7a}$, $SO_2NR^7R^{7a}$, and SO_2R^{7b};

R^6 is independently selected from the groups: halo, -CN, NO_2, C_{1-4} alkyl, C_{1-4} haloalkyl, NR^7R^{7a}, $NR^8NR^8R^{8a}$, $NR^7C(O)OR^7$, $NR^7C(O)R^{7b}$, =O, OR^7, COR^7, CO_2R^7, $CONR^7R^{7a}$, $NHC(O)NR^7R^{7a}$, $NHC(S)NR^7R^{7a}$, $SO_2N^7R^{7a}$, SO_2R^{7b}, C_{3-10} membered carbocycle substituted with 0-5 R^4, and 5-10 membered heterocycle containing from 1-4 heteroatoms selected from O, N, and S, substituted with 0-3 R^7;

R^7 is independently selected from the groups: H, halo, -CN, NO_2, C_{1-4} haloalkyl, $R^8R^{8a}N(CR^9R^{9a})m$, $NR^8NR^8R^{8a}$, $NR^8C(O)OR^8$, $NR^8C(O)R^8$, =O, $R^8O(CR^9R^{9a})m$, COR^8, CO_2R^8, $CONR^8R^{8a}$, $NHC(O)NR^8R^{8a}$, $NHC(S)NR^8R^{8a}$, $SO_2NR^8R^{8a}$, SO_2R^{8b}, C_{1-4}alkyl, C_{3-6}cycloalkyl, C_{4-10}cycloalkylalkyl, phenyl, and benzyl;

R^{7a} is independently selected from the groups: H, C_{1-4} alkyl, C_{3-6} cycloalkyl, C_{4-10} cycloalkylalkyl, phenyl, and benzyl;

alternatively, R^7 and R^{7a}, together with the atoms to which they are attached, form a heterocycle having 4-8 atoms in the ring and containing an additional 0-1 N, S, or O atom and substituted with 0-3 R^{7c};—

R^{7b} is independently selected from the groups: H, C_{1-4} alkyl, C_{3-6} cycloalkyl, C_{4-10} cycloalkylalkyl, phenyl, and benzyl;

R^{7c} is independently selected from the groups: halo, -CN , N_3, NO_2, C_{1-4} alkyl, C_{3-6} cycloalkyl, C_{4-10} cycloalkylalkyl, C_{1-4} haloalkyl, NR^7R^{7b}, $R^8R^{8a}N(CR^9R^{9a})m$, =O, OR^7, $R^8O(CR^9R^{9a})m$, COR^7, CO_2R^7, $CONR^7R^{7b}$, $NHC(O)NR^7R^{7b}$, $NHC(S)NR^7R^{7b}$, $NR^7C(O)OR^{7b}$, $NR^7C(O)R^{7b}$, $C(=NR^8)R^{8a}$, $C(=NR^8)NR^{8a}R^{8b}$, $SO_2NR^7R^{7b}$, SO_2R^{7b}, and 5-10 membered heterocycle containing from 1-4 heteroatoms selected from O, N, and S;

R^8 is independently selected from the groups: H, C_{1-4} alkyl, C_{3-6} cycloalkyl, C_{4-10} cycloalkylalkyl, phenyl and benzyl;

R^{8a} is independently selected from the groups: H, C_{1-4} alkyl, C_{3-6} cycloalkyl, C_{4-10} cycloalkylalkyl, phenyl and benzyl;

alternatively, R^8 and R^{8a}, together with the atoms to which they are attached, form a heterocycle having 4-8 atoms in the ring and containing an additional 0-1 N, S, or O atom;

R^{8b} is independently selected from the groups: H, C_{1-4} alkyl, C_{3-6} cycloalkyl, C_{4-10} cycloalkylalkyl, phenyl and benzyl;

R^9 is idependently selected from the groups: H, C_{1-4} alkyl;

R^{9a} is independently selected from the groups: H, C_{1-4} alkyl;

R^{10} is independently selected from the groups: NR^7R^{7a}, C_{3-10} membered carbocycle substituted with 0-3 R^7, and 5-10 membered heterocycle containing from 1-4 heteroatoms selected from O, N, and S, substituted with 0-3 R^7; and

m is independently selected from 0, 1, 2, 3, and 4;

or a pharmaceutically acceptable salt thereof, an N-oxide form thereof, or a stereoisomer thereof.

Claim 2 (original): A compound according to claim 1, wherein:

X is O;

R^1 is selected from the groups: C_5-C_6 membered carbocycle substituted with 0-5 R^4, and 5-6 membered heterocycle substituted with 0-5 R^5.

Claim 3 (original): A compound according to claim 1, wherein:

X is O;

R^1 is a C_5-C_6 membered carbocycle substituted with 0-5 R^4, wherein the carbocycle is an aryl,cycloalkyl, or cycloalkenyl group.

Claim 4 (original): A compound according to claim 1, wherein:

X is O;

R^1 is phenyl substituted with 0-5 R^4.

Claim 5 (original): A compound according to claim 1, wherein:

X is O;

R^1 is a C_5-C_6 membered cycloalkyl group substituted with 0-5 R^4, wherein the cycloalkyl is cyclohexyl, cyclopentyl.

Claim 6 (original): A compound according to claim 1, wherein:

X is O;

R^1 is a C_5-C_6 membered cycloalkenyl group substituted with 0-5 R^4, wherein the cycloalkenyl group is cyclohexenyl, cyclopentenyl.

Claim 7 (original): A compound according to claim 1, wherein:

X is O;

R^1 is a C_5-C_7 membered heterocycle substituted with 0-5 R^5, wherein the heterocycle is a heteroaryl,heterocyclenyl, or heterocyclyl group.

Claim 8 (original): A compound according to claim 1, wherein:

X is O;

R^1 is a C_5-C_6 membered heteroaryl substituted with 0-5 R^5, wherein the heteroaryl is pyrazinyl, thienyl, isothiazolyl, oxazolyl, pyrazolyl, furazanyl, pyrrolyl, 1,2,4-thiadiazolyl, pyridazinyl, quinoxalinyl, phthalazinyl, imidazo[1,2-a]pyridine, imidazo[2,1-b]thiazolyl, benzofurazanyl, azaindolyl, benzimidazolyl, benzothienyl, thienopyridyl, thienopyrimidyl, pyrrolopyridyl, imidazopyridyl, benzoazaindole, 1,2,4-triazinyl, benzthiazolyl, furanyl, imidazolyl, indolyl, indolizinyl, isoxazolyl, isoquinolinyl, isothiazolyl, oxadiazolyl, pyrazinyl, pyridazinyl, pyrazolyl, pyridyl, pyrimidinyl, pyrrolyl, quinazolinyl, quinolinyl, 1,3,4-thiadiazolyl, thiazolyl, thienyl or triazolyl.

Claim 9 (original): A compound according to claim 1, wherein:

X is O;

R^1 is a C_5-C_6 membered heteroaryl substituted with 0-5 R^5, wherein the heteroaryl is pyrazinyl, pyridazinyl, pyridyl, pyrimidinyl, thiazolyl or thienyl.

Claim 10 (original): A compound according to claim 1, wherein:

X is O;

R^1 is a C_5-C_6 membered heterocyclyl substituted with 0-5 R^5, wherein the heterocyclyl is tetrahydropyranyl, pyrrolidinyl, imidazolidinyl, pyrazolidinyl, piperidinyl, morpholinyl, thiomorpholinyl, or piperazinyl.

Claim 11 (original): A compound according to claim 1, wherein:

X is O;

R^1 is a C_5-C_6 membered heterocyclyl substituted with 0-5 R^5, wherein the heterocyclyl is tetrahydropyranyl or morpholinyl.

Claim 12 (original): A compound according to claim 1, wherein:

X is O;

R^1 is a C_5-C_6 membered heterocyclenyl group substituted with 0-5 R^5, wherein the heterocyclenyl group is 1,2,3,4- tetrahydrohydropyridine, 1,2-dihydropyridyl, 1,4-dihydropyridyl, 1,2,3,6-tetrahydropyridine, 1,4,5,6-tetrahydropyrimidine, 2-pyrrolinyl, 3-pyrrolinyl, 2-imidazolinyl, 2-pyrazolinyl, 3,4-dihydro-2*H*-pyran, or dihydrofuranyl.

Claim 13 (original): A compound according to claim 1, wherein:

X is O;

R^3 is selected from the groups: H, C_{1-4} alkyl.

Claim 14 (original): A compound according to claim 1, wherein:

X is O;

R^3 is methyl.

Claim 15 (original): A compound according to claim 1, wherein:

X is O;

R^2 is a C_{3-10} membered carbocycle substituted with 0-5 R^4, or a 3-10 membered heterocycle containing from 1-4 heteroatoms selected from O, N, and S and substituted with 0-5 R^5.

Claim 16 (original): A compound according to claim 1, wherein:

X is O;

R^2 is C_5-C_6 membered carbocycle substituted with 0-5 R^4, wherein the carbocycle is an aryl,cycloalkyl, or cycloalkenyl group.

Claim 17 (original): A compound according to claim 1, wherein:

X is O;

R^2 is phenyl substituted with 0-5 R^4.

Claim 18 (original): A compound according to claim 1, wherein:

X is O;

R^2 is cycloalkyl substituted with 0-5 R^4, a C_5-C_6 membered cycloalkyl group substituted with 0-5 R^4, wherein the cycloalkyl is cyclohexyl, cyclopentyl.

Claim 19 (original): A compound according to claim 1, wherein:

X is O;

R^2 is a C_5-C_6 membered cycloalkenyl group substituted with 0-5 R^4, wherein the cycloalkenyl group is cyclohexenyl, cyclopentenyl.

Clain 20 (original): A compound according to claim 1, wherein:

X is O;

R^2 is a C_5-C_7 membered heterocycle substituted with 0-5 R^5, wherein the heterocycle is a heteroaryl,heterocyclenyl, or heterocyclyl group.

Claim 21 (original): A compound according to claim 1, wherein:

X is O;

R^2 is a C_5-C_6 membered heteroaryl substituted with 0-5 R^5, wherein the heteroaryl is pyrazinyl, thienyl, isothiazolyl, oxazolyl, pyrazolyl, furazanyl, pyrrolyl, 1,2,4-thiadiazolyl, pyridazinyl, quinoxalinyl, phthalazinyl, imidazo[1,2-a]pyridine, imidazo[2,1-b]thiazolyl, benzofurazanyl, azaindolyl, benzimidazolyl, benzothienyl, thienopyridyl, thienopyrimidyl, pyrrolopyridyl, imidazopyridyl, benzoazaindole, 1,2,4-triazinyl, benzthiazolyl, furanyl, imidazolyl, indolyl, indolizinyl, isoxazolyl, isoquinolinyl, isothiazolyl, oxadiazolyl, pyrazinyl, pyridazinyl, pyrazolyl, pyridyl, pyrimidinyl, pyrrolyl, quinazolinyl, quinolinyl, 1,3,4-thiadiazolyl, thiazolyl, thienyl or triazolyl.

Claim 22 (original): A compound according to claim 1, wherein:

X is O;

R^2 is a C_5-C_6 membered heteroaryl substituted with 0-5 R^5, wherein the heteroaryl is pyrazinyl, pyridazinyl, pyridyl, pyrimidinyl, thiazolyl or thienyl.

Claim 23 (orignal): A compound according to claim 1, wherein:

X is O;

R^2 is a C_5-C_6 membered heterocyclyl substituted with 0-5 R^5, wherein the heterocyclyl is tetrahydropyranyl, pyrrolidinyl, imidazolidinyl, pyrazolidinyl, piperidinyl, morpholinyl, thiomorpholinyl, or piperazinyl.

Claim 24 (original): A compound according to claim 1, wherein:

X is O;

R^2 is a C_5-C_6 membered heterocyclenyl group substituted with 0-5 R^5, wherein the heterocyclenyl group is 1,2,3,4- tetrahydrohydropyridine, 1,2-dihydropyridyl, 1,4-dihydropyridyl, 1,2,3,6-tetrahydropyridine, 1,4,5,6-tetrahydropyrimidine, 2-pyrrolinyl, 3-pyrrolinyl, 2-imidazolinyl, 2-pyrazolinyl, 3,4-dihydro-2*H*-pyran, or dihydrofuranyl.

Claim 25 (original): A compound according to claim 1, wherein:

X is O;

R^2 is phenyl substituted with 1-5 R^4.

Claim 26 (original): A compound according to claim 1, wherein:

X is O;

R^2 is phenyl substituted with 1-4 R^4.

Claim 27 (original): A compound according to claim 1, wherein:

X is O;

R^2 is phenyl substituted with 1-3 R^4.

Claim 28 (original): A compound according to claim 1, wherein:

X is O;

R^2 is phenyl substituted with 1-2 R^4.

Claim 29 (original): A compound according to claim 1, wherein:

X is O;

R^2 is phenyl substituted with R^4;

R^4 is a 5-10 membered heterocycle containing from 1-4 heteroatoms selected from O, N, and S, wherein the heterocycle is a heteroaryl,heterocyclenyl, or heterocyclyl group.

Claim 30 (original): A compound according to claim 1, wherein:

X is O;

R^2 is phenyl substituted with R^4;

R^4 is a 5-6 membered heteroaryl containing from 1-4 heteroatoms selected from O, N, and S, which is substituted with 0-5 R^5.

Claim 31 (original): A compound according to claim 1, wherein:

X is O;

R^2 is phenyl substituted with R^4;

R^4 is NR^7R^{7a}.

Claim 32 (original): A compound according to claim 1, wherein:

X is O;

R^2 is phenyl substituted with R^4;

R^4 is NR^7R^{7a};

R^7 and R^{7a}, together with the atoms to which they are attached, form a heterocycle having 4-8 atoms in the ring and containing an additional 0-1 N, S, or O atom and substituted with 0-3 R^{7c}; and

R^{7c} is independently selected from the groups: halo, -CN , N_3, NO_2, C_{1-4} alkyl, C_{3-6} cycloalkyl, C_{4-10} cycloalkylalkyl, C_{1-4} haloalkyl, NR^7R^{7b}, $R^8R^{8a}N(CR^9R^{9a})m$, =O, OR^7, $R^8O(CR^9R^{9a})m$, COR^7, CO_2R^7, $CONR^7R^{7b}$, $NHC(O)NR^7R^{7b}$, $NHC(S)NR^7R^{7b}$, $NR^7C(O)OR^{7b}$, $NR^7C(O)R^{7b}$, $C(=NR^8)R^{8a}$, $C(=NR^8)NR^{8a}R^{8b}$, $SO_2NR^7R^{7b}$, SO_2R^{7b}, and 5-10 membered heterocycle containing from 1-4 heteroatoms selected from O, N, and S.

Claim 33 (original): A compound according to claim 1, wherein:

X is O;

R^2 is phenyl substituted with R^4;

R^4 is NR^7R^{7a};

R^7 and R^{7a}, together with the atoms to which they are attached, form a heterocycle having 6-7 atoms in the ring and containing an additional 0-1 N atoms and substituted with 0-3 R^{7c}; and

R^{7c} is independently selected from the groups: halo, -CN , N_3, NO_2, C_{1-4} alkyl, C_{3-6} cycloalkyl, C_{4-10} cycloalkylalkyl, C_{1-4} haloalkyl, NR^7R^{7b}, $R^8R^{8a}N(CR^9R^{9a})m$, =O, OR^7, $R^8O(CR^9R^{9a})m$, COR^7, CO_2R^7, $CONR^7R^{7b}$, $NHC(O)NR^7R^{7b}$, $NHC(S)NR^7R^{7b}$, $NR^7C(O)OR^{7b}$, $NR^7C(O)R^{7b}$, $C(=NR^8)R^{8a}$, $C(=NR^8)NR^{8a}R^{8b}$, $SO_2NR^7R^{7b}$, SO_2R^{7b}, and 5-10 membered heterocycle containing from 1-4 heteroatoms selected from O, N, and S.

Claim 34 (original): A compound according to claim 1, wherein:

X is O;

R^2 is phenyl substituted with R^4;

R^4 is NR^7R^{7a};

R^7 and R^{7a}, together with the atoms to which they are attached, form a 6-7 membered heterocyclyl group or a 6-7 membered heterocyclenyl group, substituted with 0-3 R^{7c}; and

R^{7c} is independently selected from the groups: halo, -CN , N_3, NO_2, C_{1-4} alkyl, C_{3-6} cycloalkyl, C_{4-10} cycloalkylalkyl, C_{1-4} haloalkyl, NR^7R^{7b}, $R^8R^{8a}N(CR^9R^{9a})m$, =O, OR^7, $R^8O(CR^9R^{9a})m$, COR^7, CO_2R^7, $CONR^7R^{7b}$, $NHC(O)NR^7R^{7b}$, $NHC(S)NR^7R^{7b}$, $NR^7C(O)OR^{7b}$, $NR^7C(O)R^{7b}$, $C(=NR^8)R^{8a}$, $C(=NR^8)NR^{8a}R^{8b}$, $SO_2NR^7R^{7b}$, SO_2R^{7b}, and 5-10 membered heterocycle containing from 1-4 heteroatoms selected from O, N, and S.

Claim 35 (original): A compound according to claim 1, wherein:

X is O;

R^2 is phenyl substituted with R^4;

R^4 is NR^7R^{7a};

R^7 and R^{7a}, together with the atoms to which they are attached, form a 6-7 membered heterocyclyl group substituted with 0-3 R^{7c}, wherein the heterocyclyl group is piperazinyl, or homopiperazinyl, and

R^{7c} is independently selected from the groups: halo, -CN , N_3, NO_2, C_{1-4} alkyl, C_{3-6} cycloalkyl, C_{4-10} cycloalkylalkyl, C_{1-4} haloalkyl, NR^7R^{7b}, $R^8R^{8a}N(CR^9R^{9a})m$, =O, OR^7, $R^8O(CR^9R^{9a})m$, COR^7, CO_2R^7, $CONR^7R^{7b}$, $NHC(O)NR^7R^{7b}$, $NHC(S)NR^7R^{7b}$, $NR^7C(O)OR^{7b}$, $NR^7C(O)R^{7b}$, $C(=NR^8)R^{8a}$, $C(=NR^8)NR^{8a}R^{8b}$, $SO_2NR^7R^{7b}$, SO_2R^{7b}, and 5-10 membered heterocycle containing from 1-4 heteroatoms selected from O, N, and S.

Claim 36 (original): A compound according to claim 1, wherein:

X is O;

R^2 is phenyl substituted with R^4;

R^4 is NR^7R^{7a};

R^7 and R^{7a}, together with the atoms to which they are attached, form a 6-7 membered heterocyclenyl group substituted with 0-3 R^{7c}, wherein the heterocyclenyl group is ,2,3,4-tetrahydrohydropyridine, 1,2-dihydropyridyl, 1,4-dihydropyridyl, 1,2,3,6-tetrahydropyridine, or 1,4,5,6-tetrahydropyrimidine; and

R^{7c} is independently selected from the groups: halo, -CN , N_3, NO_2, C_{1-4} alkyl, C_{3-6} cycloalkyl, C_{4-10} cycloalkylalkyl, C_{1-4} haloalkyl, NR^7R^{7b}, $R^8R^{8a}N(CR^9R^{9a})m$, =O, OR^7, $R^8O(CR^9R^{9a})m$, COR^7, CO_2R^7, $CONR^7R^{7b}$, $NHC(O)NR^7R^{7b}$, $NHC(S)NR^7R^{7b}$, $NR^7C(O)OR^{7b}$, $NR^7C(O)R^{7b}$, $C(=NR^8)R^{8a}$, $C(=NR^8)NR^{8a}R^{8b}$, $SO_2NR^7R^{7b}$, SO_2R^{7b}, and 5-10 membered heterocycle containing from 1-4 heteroatoms selected from O, N, and S.

Claim 37 (original): A compound according to claim 1, wherein:

R^{7c} is independently selected from the groups: C_{1-4} alkyl, C_{3-6} cycloalkyl, C_{4-10} cycloalkylalkyl, $NR^{7}R^{7b}$, and 5-10 membered heterocycle containing from 1-4 heteroatoms selected from O, N, and S.

Claim 38 (original): A compound according to claim 1, wherein the compound is selected from:

3-(4-piperazinophenyl)-5-((N-methyl- N-(2-pyridinyl)amino)carbamoylamino)indeno[1,2-c]pyrazol-4-one;

3-(4-(4-methylpiperazino)phenyl)-5-((N-methyl- N-(2-pyridinyl)amino)carbamoylamino)indeno[1,2-c]pyrazol-4-one;

3-(4-homopiperazinophenyl)-5-((N-methyl- N-(2-pyridinyl)amino)carbamoylamino)indeno[1,2-c]pyrazol-4-one;

3-(4-(4-methylhomopiperazino)phenyl)-5-((N-methyl- N-(2-pyridinyl)amino)carbamoylamino)indeno[1,2-c]pyrazol-4-one;

3-(4-piperazinophenyl)-5-((N-methyl-N-(4-pyridinyl)amino)carbamoylamino)indeno[1,2-c]pyrazol-4-one;

3-(4-piperazinophenyl)-5-((N-methyl-N-(2-pyrazinyl)amino)carbamoylamino)indeno[1,2-c]pyrazol-4-one;

3-(4-piperazinophenyl)-5-((N-methyl-N-(2-pyrimidinyl)amino)carbamoylamino)indeno[1,2-c]pyrazol-4-one;

3-(4-piperazinophenyl)-5-((N-methyl-N-(2-thiazolyl)amino)carbamoylamino)indeno[1,2-c]pyrazol-4-one;

3-(4-piperazinophenyl)-5-((N-methyl-N-(3-pyridinyl)amino)carbamoylamino)indeno[1,2-c]pyrazol-4-one;

3-(4-(4-methylpiperazino)phenyl)-5-((N-methyl-N-(2-pyrazinyl)amino)carbamoylamino)indeno[1,2-c]pyrazol-4-one;

3-(4-(4-methylpiperazino)phenyl)-5-((N-methyl-N-(2-thiazolyl)amino)carbamoylamino)indeno[1,2-c]pyrazol-4-one;

3-(4-(4-methylpiperazino)phenyl)-5-((N-methyl-N-(3-pyridinyl)amino)carbamoylamino)indeno[1,2-c]pyrazol-4-one;

3-(4-piperazinophenyl)-5-((N-methyl-N-(4-tetrahydropyranyl)amino)carbamoylamino)indeno[1,2-c]pyrazol-4-one;

3-(4-(4-methylpiperazino)phenyl)-5-((N-methyl- N-(4-tetrahydropyranyl)amino)carbamoylamino)-indeno[1,2-c]pyrazol-4-one;

3-(4-(4-ethylpiperazino)phenyl)-5-((N-methyl- N-(4-tetrahydropyranyl)amino)carbamoylamino)indeno[1,2-c]pyrazol-4-one;

3-(4-(4-isopropylpiperazino)phenyl)-5-((N-methyl- N-(4-tetrahydropyranyl)amino)carbamoylamino)-indeno[1,2-c]pyrazol-4-one;

3-(4-(4-piperazinophenyl)-5-((N-methyl-N-cyclohexylamino)carbamoylamino)-indeno[1,2-c]pyrazol-4-one;

3-(4-(4-methylpiperazino)phenyl)-5-((N-methyl-N-cyclohexylamino)carbamoylamino)-indeno[1,2-c]pyrazol-4-one;

3-(4-(4-ethylpiperazino)phenyl)-5-((N-methyl-N-cyclohexylamino)carbamoylamino)indeno[1,2-c]pyrazol-4-one;

3-(4-(4-isopropylpiperazino)phenyl)-5-((N-methyl-N-cyclohexylamino)carbamoylamino)-indeno[1,2-c]pyrazol-4-one;

3-(4-piperazinophenyl)-5-((N-methyl-N-(1-methylpiperidin-4-yl)amino)carbamoylamino)indeno[1,2-c]pyrazol-4-one;

3-(4-homopiperazinophenyl)-5-((N-methyl-N-(4-tetrahydropyranyl)amino)carbamoylamino)indeno[1,2-c]pyrazol-4-one;

3-(4-(4-methylhomopiperazino)phenyl)-5-((N-methyl-N-(4-tetrahydropyranyl)amino)carbamoylamino)-indeno[1,2-c]pyrazol-4-one;

3-(4-(4-ethylhomopiperazino)phenyl)-5-((N-methyl-N-(4-tetrahydropyranyl)amino)carbamoylamino)-indeno[1,2-c]pyrazol-4-one;

3-(4-(4-isopropylhomopiperazino)phenyl)-5-((N-methyl-N-(4-tetrahydropyranyl)amino)carbamoylamino)-indeno[1,2-c]pyrazol-4-one;

3-(4-(4-(N,N-dimethylamino)piperidino)phenyl)-5-((N-methyl-N-(4-tetrahydropyranyl)amino)carbamoylamino)-indeno[1,2-c]pyrazol-4-one;

3-(4-(4-pyrrolidinopiperidino)phenyl)-5-((N-methyl-N-(4-tetrahydropyranyl)amino)carbamoylamino)-indeno[1,2-c]pyrazol-4-one;

3-(4-(4-piperidinopiperidino)phenyl)-5-((N-methyl-N-(4-tetrahydropyranyl)amino)carbamoylamino)-indeno[1,2-c]pyrazol-4-one;

3-(2,4-dimethylthiazol-5-yl)-5-((N-methyl-N-(4-tetrahydropyranyl)amino)carbamoylamino)indeno[1,2-c]pyrazol-4-one;
or pharmaceutically acceptable salt form thereof.

Claims 39-61 (cancelled)